

# **ORNL Cross-Section Processing Status**

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# AMPX Status

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- ENDF/B-VI Multigroup Library Development
  - Corrected processing problems with collision kinematics treatment
  - Corrected various systematic errors in processing procedures
  - Generated 238-group ENDF/B-VI library for testing
  - Currently performing various benchmark calculations with the library in the SCALE 5 code system
  - Good results obtained for various benchmark calculations; however, some results are consistently 1-2% below critical.
  - Currently investigating multigroup scattering matrices to resolve discrepancies with benchmark results

# AMPX Status

- AMPX developments to support continuous-energy version of KENO V.a
  - Developed new processing modules to support Point KENO development effort
  - Point KENO cross-section library components
    - $\bar{\nu}(E)$  delayed and prompt
    - 1-D continuous-energy cross sections as a function of temperature  $\sigma(E,T)$
    - 2-D joint probability distributions for particle collisions  $f(E \rightarrow E', \mu)$
    - Probability tables for unresolved resonance region
  - MONTEGO and JAMAICAN modules
    - Generate joint probability density functions (PDF) and cumulative distribution functions (CDF)
    - Produce tabular PDFs and CDFs in lab system
  - PLATINUM module developed to assemble Point KENO cross-section file for a nuclide
- Used AMPX to generate test ENDF/B-VI library for Point KENO
  - 50 Nuclides

# Point KENO V.a Testing

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- Prototypic Point KENO cross-section data generated for 50 ENDF/B-VI Release 7 nuclides

$^{27}\text{Al}$	$^{52}\text{Cr}$	$^{56}\text{Fe}$	$^{55}\text{Mn}$	$^{31}\text{P}$	$^{242}\text{Pu}$	$^{30}\text{Si}$	$^{238}\text{U}$
$^{241}\text{Am}$	$^{53}\text{Cr}$	$^{57}\text{Fe}$	$\text{Mo}$	$^{236}\text{Pu}$	$^{243}\text{Pu}$	$^{232}\text{U}$	
$^{10}\text{B}$	$\text{C}$	$^{58}\text{Fe}$	$^{14}\text{N}$	$^{237}\text{Pu}$	$^{244}\text{Pu}$	$^{233}\text{U}$	
$^{11}\text{B}$	$^{63}\text{Cu}$	$^1\text{H}$	$^{23}\text{Na}$	$^{238}\text{Pu}$	$\text{S}$	$^{234}\text{U}$	
$\text{Ca}$	$^{65}\text{Cu}$	$\text{H}_2\text{O}$	$^{58}\text{Ni}$	$^{239}\text{Pu}$	$\text{Si}$	$^{235}\text{U}$	
$\text{Cd}$	$^{19}\text{F}$	$\text{CH}_2$	$^{60}\text{Ni}$	$^{240}\text{Pu}$	$^{28}\text{Si}$	$^{236}\text{U}$	
$^{50}\text{Cr}$	$^{54}\text{Fe}$	$\text{K}$	$^{16}\text{O}$	$^{241}\text{Pu}$	$^{29}\text{Si}$	$^{237}\text{U}$	



# Point KENO V.a Testing

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- Point KENO V.a used to calculate 54 different criticality test problems
  - 33 KENO test problems
    - Uranium metal (single units and arrays)
    - Uranyl fluoride and nitrate solutions
    - Mixed metal and solution problems
    - Various geometrical and reflector configurations
  - 21 benchmark problems
    - LWR-type  $\text{UO}_2$  fuel pin lattices
    - Green-block experiments: homogenized U in paraffin blocks
    - Uranyl fluoride and nitrate solutions (low and high enriched)
- Calculations performed on DEC Alpha XP1000 workstation

# Point KENO V.a Testing

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Case ID	Description	Point KENO ENDF/B-VI Rel. 7	KENO V.a 199-group ENDF/B-VI Rel. 3	KENO V.a 238-group ENDF/B-V
cas01	U(2.35)O <sub>2</sub> pin lattice	0.9950 ± 0.0016	0.9947 ± 0.0015	0.9937 ± 0.0015
cas07	U(2.35)O <sub>2</sub> pin lattice	0.9965 ± 0.0018	0.9934 ± 0.0016	0.9956 ± 0.0015
cas19	U(2.35)O <sub>2</sub> pin lattice	1.0008 ± 0.0016	0.9937 ± 0.0022	0.9903 ± 0.0014
cas34	U(2.46)O <sub>2</sub> pin lattice	0.9997 ± 0.0016	0.9973 ± 0.0011	0.9925 ± 0.0015
cas82	U(2.0) in Paraffin	1.0003 ± 0.0014	1.0241 ± 0.0015	1.0021 ± 0.0016
cas83	U(2.0) in Paraffin	1.0011 ± 0.0019	1.0194 ± 0.0017	0.9992 ± 0.0017
cas84	U(3.0) in Paraffin	1.0129 ± 0.0021	1.0357 ± 0.0017	1.0118 ± 0.0017
cas85	U(3.0) in Paraffin	1.0177 ± 0.0016	1.0407 ± 0.0020	1.0123 ± 0.0020
cas88	U(93.2)O <sub>2</sub> (NO <sub>3</sub> ) <sub>2</sub>	1.0026 ± 0.0023	1.0029 ± 0.0025	1.0090 ± 0.0021